This listing of claims will replace all prior versions, and listings, of claims in the application:

Listing of Claims:

1. (Previously Presented) A compound of Formula I:

A - D - B (I)

or a pharmaceutically acceptable salt thereof, wherein

D is -NH-C(O)-NH-,

A is a substituted moiety of the formula:

-L-M-L^1,

wherein L is phenyl, optionally substituted by halogen, up to per-halo, and Wn, where n is 0-3;

wherein each W is independently selected from the group consisting of C_1-C_5 linear or branched alkyl, C_1-C_5 linear or branched haloalkyl up to perhaloalkyl and C_1-C_3 alkoxy L^1 is selected from pyridinyl substituted by -C(O)R_x, and

optionally substituted with 1-3 additional substituents independently selected from the group consisting of R^7 and halogen;

wherein R_x is NR_aR_b and R_a and R_b are

A) independently

a) hydrogen,

b) C_1-C_{10} alkyl,

c) C_6 aryl,

d) pyridinyl

e) substituted C_{1-10} alkyl,

f) substituted C_6 aryl,

g) substituted pyridinyl

h) -phenylpiperazine(pyridinyl),

i) -phenylmorpholinyl,

j) -ethylmorpholinyl,

k) -ethylpiperidyl,

l) -methyl pyrrolidinyl,

m) -methyl tetrahydrofuryl,

or

n) $-C_2H_4NH$(phenyl);

where when R_a and R_b are a substituted group, they are substituted by

a) halogen up to per halo,

b) hydroxy,

c) $-N(CH_3)_2$,

d) C_1-C_{10} alkyl,

e) C_1-C_{10} alkoxy,

f) halosubstituted C_{1-6} alkyl, or

g) $-OSi(Pr\text{-}i)_{3;}$ or

B) R_a and R_b together form piperazine or a substituted piperazine with substituents selected from the group consisting of

a) halogen,

b) hydroxy,

c) C_{1-10} alkyl,

d) pyridinyl

e) C_{1-10} alkoxy,

f) C_6 aryl,

g) halo substituted C_6 aryl, and

h) N-(4-acetylphenyl);

M is selected from the group consisting of oxygen and sulfur;

and

B is

phenyl, substituted with 1-3 substituents independently selected from the group consisting of halogen and R^7,

and R^7 is

(a) C_1-C_6 linear or branched alkyl, optionally substituted with 1-3 halogen substituents; or

(b) C_1-C_6 linear or branched alkoxy.

2. (Canceled)

3. (Previously Presented) A compound as in claim 1 wherein M is oxygen .

4. (Previously Presented) A compound as in claim 1 wherein the cyclic structures of B and L bound directly to D are substituted in the ortho position by hydrogen.

5. (Canceled)

6. (Previously Presented) A compound of claim 1 wherein B of Formula I is phenyl, substituted with 1-3 substituents independently selected from the group consisting of chlorine, C_1-C_6 alkoxy or up to per halo substituted C_1-C_6 alkyl.

7. (Previously Presented) A compound of claim 3 wherein B of Formula I is phenyl, substituted with 1-3 substituents independently selected from the group consisting of chlorine, C_1-C_6 alkoxy, or substituted C_1-C_6 alkyl, substituted by one or more halogen substituents.

8. (Previously Presented) A compound of claim 4 wherein B of Formula I is phenyl, substituted 1 to 3 times by 1 or more substituents selected from the group consisting of- chlorine, C_1-C_6 alkoxy or up to per halo substituted C_1-C_6 alkyl.

9. (Previously Presented) A compound of claim 1, wherein L is phenyl, optionally substituted by halogen up to perhalo.

10. (Previously Presented) A compound of claim 1, wherein L is phenyl, optionally substituted with 1-3 substituents independently selected from the group consisting of halogen and C_1-C_3 alkoxy.

11. (Canceled)

12. (Canceled)

13. (Canceled)

14. (Canceled)

15. (Canceled)

16. (Canceled)

17. (Canceled)

18. (Previously Presented) A compound of claim 4, wherein M is -O- .

19. (Previously Presented) A compound of claim 8 wherein M is -O-.

20. (Previously Presented) A compound of claim 9 wherein M is -O-.

21. (Previously Presented) A compound of claim 10 wherein M is -O- .

22. (Previously Presented) A compound of claim 1 wherein L^1 is additionally substituted 1 to 3 times by one or more substituents selected from the group consisting of C_1-C_6 alkyl, halogen and C_1-C_6 alkoxy.

23. (Previously Presented) A compound of claim 3 wherein L^1 is additionally substituted 1 to 3 times by one or more substituents selected from the group consisting of C_1-C_6 alkyl, halogen and C_1-C_6 alkoxy.

24. (Previously Presented) A compound of claim 18 wherein L^1 is additionally substituted 1 to 3 times by one or more substituents selected from the group consisting of C_1-C_6 alkyl, halogen and C_1-C_6 alkoxy.

25. (Previously Presented) A compound of claim 19 wherein L^1 is additionally substituted 1 to 3 times by one or more substituents selected from the group consisting of C_1-C_6 alkyl, halogen and C_1-C_6 alkoxy.

26. (Previously Presented) A compound of claim 20 wherein L^1 is additionally substituted 1 to 3 times by one or more substituents selected from the group consisting of C_1-C_6 alkyl, halogen and C_1-C_6 alkoxy.

27. (Previously Presented) A compound of claim 21 wherein L^1 is additionally substituted 1 to 3 times by one or more substituents selected from the group consisting of C_1-C_6 alkyl, halogen and C_1-C_6 alkoxy.

28. (Canceled)

29. (Canceled)

30. (Canceled)

31. (Canceled)

32. (Canceled)

33. (Previously Presented) A compound of claim 3 wherein R_a and R_b are independently hydrogen or C_1-C_6 alkyl.

34. (Previously Presented) A compound of claim 18 wherein R_a and R_b are independently hydrogen or C_1-C_6 alkyl.

35. (Previously Presented) A compound of claim 19 wherein R_a and R_b are independently hydrogen or C_1-C_6 alkyl.

36. (Previously Presented) A compound of claim 20 wherein R_a and R_b are independently hydrogen or C_1-C_6 alkyl.

37. (Previously Presented) A compound of claim 21 wherein R_a and R_b are independently hydrogen or C_1-C_6 alkyl.

38. (Previously Presented) A compound of Formula I:

A - D - B (I)

or a pharmaceutically acceptable salt thereof, wherein

D is -NH-C(O)-NH-,

A is of the formula: -L-M-L^1, wherein

L is phenyl, optionally substituted with 1-3 substituents independently selected from the group consisting of C_1-C_5 linear or branched alkyl, C_1-C_5 linear or branched haloalkyl up to perhalo, C_1-C_3 alkoxy and halogen;

L^1 is pyridinyl, substituted by -C(O)R_x;

wherein R_x is NR_aR_b and R_a and R_b are independently

hydrogen,

C_1-C_{10} alkyl,

C_6 aryl,

pyridinyl, substituted $C_{1\text{-}10}$ alkyl,

substituted C_6 aryl, or

substituted pyridinyl,

where R_a and R_b are a substituted group, they are substituted by halogen up to per halo; and

M is selected from the group consisting of oxygen and sulfur

and

B is phenyl, substituted with 1-3 substituents independently selected from the group consisting of R^7 and halogen;

and R^7 is

(a) C_1-C_6 linear or branched alkyl, optionally substituted with 1-3 halogen substituents; or

(b) C_1-C_6 linear or branched alkoxy.

39. (Previously Presented) A compound of Formula I:

A - D - B (I)

or a pharmaceutically acceptable salt thereof, wherein

D is -NH-C(O)-NH-,

A is of the formula: -L-M-L^1,

L is phenyl,

M is -O-,

L^1 is pyridinyl substituted by -C(O)R_x ,

wherein R_x is NR_aR_b and R_a and R_b are independently

hydrogen,

C_1-C_{10} alkyl,

C_6 aryl,

pyridinyl,

substituted C_{1-10} alkyl,

substituted C_6 aryl, or

substituted pyridinyl

where R_a and R_b are a substituted group, they are substituted by halogen up to per halo,

and

B is a phenyl group substituted by trifluoromethyl or tert-butyl, and optionally additional substituents selected from the group consisting of halogen up to per halo, and W_n where n is 0-3, and each W is independently selected from the group consisting of

C_1-C_{10} alkyl,

C_1-C_{10} alkoxy,

C_6 aryl,

pyridinyl,

and substituted C_1-C_{10} alkyl, substituted by one or more substituents independently selected from the group consisting of halogen up to per halo.

40. (Previously Presented) A compound as in claim 38 wherein the cyclic structures of B and L bound directly to D are substituted in the ortho position by hydrogen.

41. (Canceled)

42. (Previously Presented) A compound as in claim 39 wherein the cyclic structures of B and L bound directly to D are substituted in the ortho position by hydrogen.

43. (Canceled)

44. (Previously Presented) A compound as in claim 38 wherein substituents for B, are selected from the group consisting of up to per halo substituted C_1-C_6 alkyl and halogen.

45. (Previously Presented) A compound as in claim 39 wherein the optional substituents for B are selected from the group consisting of up to per halo substituted C_1-C_6 alkyl and halogen.

46. (Canceled)

47. (Canceled)

48. (Canceled)

49. (Canceled)

50. (Previously Presented) A pharmaceutically acceptable salt of a compound of formula I of claim 1 which is

a) a basic salt of an organic acid or inorganic acid which is hydrochloric acid, hydrobromic acid, sulfuric acid, phosphoric acid, methanesulfonic acid, trifluoromethanesulfonic acid, benzenesulfonic acid, p-toluene sulfonic acid (tosylate salt), 1-napthalene sulfonic acid, 2-napthalene sulfonic acid, acetic acid, trifluoroacetic acid, malic acid, tartaric acid, citric acid, lactic acid, oxalic acid, succinic acid, fumaric acid, maleic acid, benzoic acid, salicylic acid, phenylacetic acid, or mandelic acid; or

b) an acid salt of an organic or inorganic base containing an alkali metal cation, an alkaline earth metal cation, an ammonium cation, an aliphatic substituted ammonium cation or an aromatic substituted ammonium cation.

51. (Canceled)

52. (Canceled)

53. (Previously Presented) A pharmaceutically acceptable salt of a compound of claim 38 which is

a) a basic salt of an organic acid or inorganic acid which is hydrochloric acid, hydrobromic acid, sulfuric acid, phosphoric acid, methanesulfonic acid, trifluoromethanesulfonic acid, benzenesulfonic acid, p-toluene sulfonic acid (tosylate salt), 1-napthalene sulfonic acid, 2-napthalene sulfonic acid, acetic acid, trifluoroacetic acid, malic acid, tartaric acid, citric acid, lactic acid, oxalic acid, succinic acid, fumaric acid, maleic acid, benzoic acid, salicylic acid, phenylacetic acid, or mandelic acid; or

b) an acid salt of an organic or inorganic base containing an alkali metal cation, an alkaline earth metal cation, an ammonium cation, an aliphatic substituted ammonium cation or an aromatic substituted ammonium cation.

54. (Previously Presented) A pharmaceutically acceptable salt of a compound of claim 39 which is

a) a basic salt of an organic acid or inorganic acid which is hydrochloric acid, hydrobromic acid, sulfuric acid, phosphoric acid, methanesulfonic acid, trifluoromethanesulfonic acid, benzenesulfonic acid, p-toluene sulfonic acid (tosylate salt), 1-napthalene sulfonic acid, 2-napthalene sulfonic acid, acetic acid, trifluoroacetic acid, malic acid, tartaric acid, citric acid, lactic acid, oxalic acid, succinic acid, fumaric acid, maleic acid, benzoic acid, salicylic acid, phenylacetic acid, or mandelic acid; or

b) an acid salt of an organic or inorganic base containing an alkali metal cation, an alkaline earth metal cation, an ammonium cation, an aliphatic substituted ammonium cation or an aromatic substituted ammonium cation.

55. (Canceled)

56. (Canceled)

57. (Canceled)

58. (Canceled)

59. (Canceled)

60. (Canceled)

61. **(Cancelled)**

62. **(Canceled)**

63. **(Canceled)**

64. **(Canceled)**

65. **(Canceled)**

66. **(Canceled)**

67. **(Canceled)**

68. **(Previously Presented)** A compound of claim 1 wherein the optional substituents on L^1 are selected from the group consisting of methyl, triflouromethyl, methoxy, Cl and F.

69. **(Previously Presented)** A compound of claim 1 wherein the substituents of B and L are independently selected from the group consisting of methyl, triflouromethyl, tert-butyl, methoxy, Cl, and F.

70. **(Canceled)**

71. **(Currently Amended)** A compound of Formula I:

A - D - B (I)

or a pharmaceutically acceptable salt thereof, wherein

D is -NH-C(O)-NH-,

A is a substituted moiety of the formula: -L-M-L^1,

wherein L is phenyl, optionally substituted with chlorine or methyl substituents ~~independently selected from the group consisting of halogen, C_1-C_5-alkyl, C_1-C_5-alkyl substituted by halogen and C_1-C_5-alkoxy~~;

L^1 is pyridinyl, substituted with -C(O)NR^aR^b ~~and optionally substituted with one or two substituents selected from the group consisting of R^7, OR^7 and halogen, wherein R^7 is hydrogen, C_1-C_5-alkyl or C_1-C_5-alkyl substituted by halogen,~~

wherein R^a and R^b independently are

a) hydrogen ~~or~~

b) ~~C_1-C_5-alkyl;~~ methyl;

c) ethyl; or

d) propyl

B is phenyl, substituted by tert-butyl or trifluoromethyl and optionally substituted with additional substituents independently selected from the group consisting of

a) halogen, or

b) methoxy ~~C_1-C_5 alkyl substituted by halogen or~~

~~c) C_1-C_4 alkoxy~~.

72. **(Canceled)**

73. **(Canceled)**

74. **(Canceled)**

75. **(Canceled)**

76. **(Canceled)**

77. **(Canceled)**

78. **(Canceled)**

79. **(Canceled)**

80. **(Canceled)**

81. **(Canceled)**

82. **(Canceled)**

83. **(Canceled)**

84. **(Canceled)**

85. **(Canceled)**

86. **(Canceled)**

87. **(Canceled)**

88. **(Canceled)**

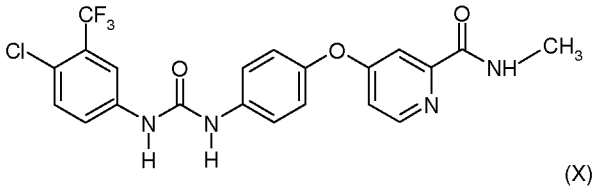
89. **(Cancelled)**

90. **(New)** A compound of claim 71 where L has no optional substituents.

91. **(New)** A compound of claim 71 where R^a is hydrogen and R^b is methyl.

92. (New) A compound of claim 71 where B is substituted by trifluoromethyl and chlorine or bromine.

93. (New) A compound which is *N*-(4-chloro-3-(trifluoromethyl) phenyl)-*N'*-(4-(2-(*N*-methylcarbamoyl)-4-pyridyloxy) phenyl) urea of the formula X



(X)

or a pharmaceutically acceptable salt thereof.

94 . (New) A compound of claim 93 which is a pharmaceutically acceptable salt of *N*-(4-chloro-3-(trifluoromethyl) phenyl)-*N'*-(4-(2-(*N*-methylcarbamoyl)-4-pyridyloxy) phenyl) urea that is a basic salt of an organic acid or an inorganic acid which is hydrochloric acid, hydrobromic acid, sulfuric acid, phosphoric acid, methanesulfonic acid, trifluoromethanesulfonic acid, benzenesulfonic acid, p-toluene sulfonic acid (tosylate salt), 1-napthalene sulfonic acid, 2-napthalene sulfonic acid, acetic acid, trifluoroacetic acid, malic acid, tartaric acid, citric acid, lactic acid, oxalic acid, succinic acid, fumaric acid, maleic acid, benzoic acid, salicylic acid, phenylacetic acid, or mandelic acid.

95. (New) A compound of claim 93 which is a tosylate salt of *N*-(4-chloro-3-(trifluoromethyl) phenyl)-*N'*-(4-(2-(*N*-methylcarbamoyl)-4-pyridyloxy) phenyl) urea.